

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011725\
 Data File : BP023631.D
 Acq On : 17 Jan 2025 16:40
 Operator : RC/JU
 Sample : Q1086-13
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AF1

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP011425.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP023631.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.317	35	39	44	rBV	125120	155209	0.94%	0.087%
2	3.723	104	108	118	rBV	463946	713729	4.33%	0.402%
3	4.146	175	180	187	rBV4	40763	62494	0.38%	0.035%
4	4.411	221	225	233	rVB2	61153	93648	0.57%	0.053%
5	5.722	442	448	462	rVB	734923	1097752	6.66%	0.619%
6	5.875	469	474	478	rBV	93782	149519	0.91%	0.084%
7	5.917	478	481	487	rVB	137395	188060	1.14%	0.106%
8	6.211	523	531	541	rBV2	40996	67789	0.41%	0.038%
9	6.446	562	571	580	rVB	97133	168426	1.02%	0.095%
10	6.534	580	586	598	rBV	103236	174753	1.06%	0.098%
11	6.787	621	629	635	rBV	44384	84131	0.51%	0.047%
12	6.969	653	660	670	rVB	676473	1109430	6.73%	0.625%
13	7.140	682	689	698	rBV	2057299	3220443	19.53%	1.815%
14	7.340	714	723	734	rBV	2407022	3711628	22.51%	2.091%
15	7.811	795	803	809	rBV	1992044	3036228	18.41%	1.711%
16	7.875	809	814	823	rVV	353555	540925	3.28%	0.305%
17	8.493	911	919	932	rVV	1998859	3022552	18.33%	1.703%
18	8.616	932	940	947	rVB	48223	89982	0.55%	0.051%
19	8.946	989	996	1005	rVV	2252852	3615557	21.92%	2.037%
20	9.063	1006	1016	1023	rVV4	83864	203122	1.23%	0.114%
21	9.416	1070	1076	1086	rVV2	91807	189216	1.15%	0.107%
22	9.511	1086	1092	1100	rVB2	57724	105857	0.64%	0.060%
23	9.663	1106	1118	1132	rBV	1332010	2448387	14.85%	1.380%
24	9.787	1133	1139	1148	rVV4	55844	130293	0.79%	0.073%
25	10.199	1202	1209	1223	rBV	3538085	5671937	34.39%	3.196%
26	10.363	1230	1237	1250	rBV	6903638	10127317	61.41%	5.706%
27	10.581	1267	1274	1286	rVB	2691777	4496456	27.27%	2.534%
28	10.710	1288	1296	1320	rBV	1876033	3383124	20.52%	1.906%
29	11.222	1377	1383	1390	rVB	96792	173198	1.05%	0.098%
30	11.393	1408	1412	1418	rVB5	33781	62453	0.38%	0.035%
31	12.163	1533	1543	1550	rBV2	55212	119871	0.73%	0.068%
32	12.334	1564	1572	1578	rBV2	35424	93085	0.56%	0.052%
33	12.410	1578	1585	1592	rVB3	55606	114720	0.70%	0.065%
34	12.604	1614	1618	1625	rBV	41197	65448	0.40%	0.037%
35	12.840	1644	1658	1671	rBV	1223818	2072230	12.57%	1.168%
36	13.269	1727	1731	1737	rVB3	49630	67051	0.41%	0.038%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP011425.MA.M
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37	13.399	1748	1753	1763	rVB3	97553	175871	1.07%	0.099%
38	13.834	1820	1827	1833	rBV	5915361	8127285	49.28%	4.579%
39	13.969	1843	1850	1856	rVB4	81249	126630	0.77%	0.071%
40	14.081	1864	1869	1870	rBV	194494	246628	1.50%	0.139%
41	14.116	1870	1875	1882	rVV	6511857	9869450	59.85%	5.561%
42	14.281	1899	1903	1910	rVB5	49146	89338	0.54%	0.050%
43	14.428	1924	1928	1932	rVV	3346950	3664881	22.22%	2.065%
44	14.463	1932	1934	1946	rVB	2599748	4748578	28.80%	2.676%
45	14.563	1946	1951	1955	rBV	676801	1069293	6.48%	0.602%
46	14.610	1955	1959	1977	rVB2	1306934	2583432	15.67%	1.456%
47	14.757	1979	1984	1992	rBV2	71426	116536	0.71%	0.066%
48	15.157	2045	2052	2061	rVV2	144683	291643	1.77%	0.164%
49	15.263	2065	2070	2075	rVB2	150680	199080	1.21%	0.112%
50	15.434	2092	2099	2110	rBV	9035903	12583072	76.30%	7.090%
51	15.540	2110	2117	2132	rVV	1499919	2123918	12.88%	1.197%
52	15.681	2136	2141	2147	rVB3	50659	91578	0.56%	0.052%
53	15.804	2156	2162	2168	rBV	1059950	1479084	8.97%	0.833%
54	15.863	2168	2172	2179	rVB8	25542	54731	0.33%	0.031%
55	16.016	2193	2198	2205	rVB3	62629	94218	0.57%	0.053%
56	16.104	2208	2213	2221	rBV6	50960	105042	0.64%	0.059%
57	16.198	2225	2229	2231	rBV2	43402	59302	0.36%	0.033%
58	16.240	2231	2236	2237	rVV	809840	868681	5.27%	0.489%
59	16.263	2237	2240	2253	rVB	1014975	1709333	10.37%	0.963%
60	16.934	2350	2354	2359	rBV3	51780	92422	0.56%	0.052%
61	17.016	2362	2368	2375	rVB3	217455	367268	2.23%	0.207%
62	17.222	2397	2403	2410	rBV2	4496136	7107023	43.10%	4.004%
63	17.322	2414	2420	2429	rVV	9802561	13927856	84.46%	7.848%
64	17.487	2444	2448	2453	rBV	102059	142378	0.86%	0.080%
65	17.540	2453	2457	2461	rVB	76694	93695	0.57%	0.053%
66	17.610	2463	2469	2477	rBV	321965	459315	2.79%	0.259%
67	17.687	2477	2482	2488	rBV	250069	371820	2.25%	0.210%
68	17.745	2488	2492	2495	rVV4	37426	77719	0.47%	0.044%
69	17.781	2495	2498	2501	rVB	52968	64218	0.39%	0.036%
70	17.828	2503	2506	2513	rVV2	55744	121834	0.74%	0.069%
71	17.892	2515	2517	2520	rVV3	45774	57681	0.35%	0.033%
72	17.934	2520	2524	2539	rVB3	156712	256727	1.56%	0.145%
73	18.057	2539	2545	2551	rBV5	35274	88981	0.54%	0.050%
74	18.175	2561	2565	2572	rBV	296495	457497	2.77%	0.258%
75	18.257	2573	2579	2587	rVB3	115474	259658	1.57%	0.146%
76	18.498	2613	2620	2627	rVB5	97629	194861	1.18%	0.110%

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77	18.639	2640	2644	2653	rVB4	76954	153056	0.93%	0.086%
78	18.945	2694	2696	2705	rVB3	92462	123256	0.75%	0.069%
79	19.045	2709	2713	2717	rBV4	30294	57624	0.35%	0.032%
80	19.157	2728	2732	2740	rBV3	72077	118259	0.72%	0.067%
81	19.245	2743	2747	2755	rVB	119266	169690	1.03%	0.096%
82	19.322	2756	2760	2766	rVB	105140	146569	0.89%	0.083%
83	19.504	2787	2791	2797	rBV	167597	255766	1.55%	0.144%
84	19.592	2801	2806	2816	rVV	523088	880884	5.34%	0.496%
85	19.704	2818	2825	2832	rVV	11199704	16490902	100.00%	9.292%
86	19.763	2832	2835	2839	rVB3	74931	108151	0.66%	0.061%
87	19.845	2846	2849	2853	rBV3	57262	75596	0.46%	0.043%
88	20.004	2872	2876	2883	rVB2	88832	139833	0.85%	0.079%
89	20.481	2953	2957	2960	rBV2	240242	298781	1.81%	0.168%
90	20.551	2964	2969	2975	rVV3	200220	334577	2.03%	0.189%
91	21.092	3057	3061	3070	rBV3	103643	186016	1.13%	0.105%
92	21.192	3074	3078	3085	rBV2	204484	413023	2.50%	0.233%
93	21.375	3106	3109	3117	rVB2	190381	278431	1.69%	0.157%
94	21.492	3126	3129	3136	rVB2	120944	157635	0.96%	0.089%
95	21.639	3151	3154	3157	rBV	184775	219098	1.33%	0.123%
96	21.692	3157	3163	3175	rVB2	4401254	7518713	45.59%	4.236%
97	22.110	3230	3234	3242	rVB2	110889	171474	1.04%	0.097%
98	23.057	3390	3395	3406	rVB	536963	1040671	6.31%	0.586%
99	24.880	3696	3705	3723	rBV	5242267	15044051	91.23%	8.477%
100	25.127	3737	3747	3764	rVB	2481162	7651692	46.40%	4.311%

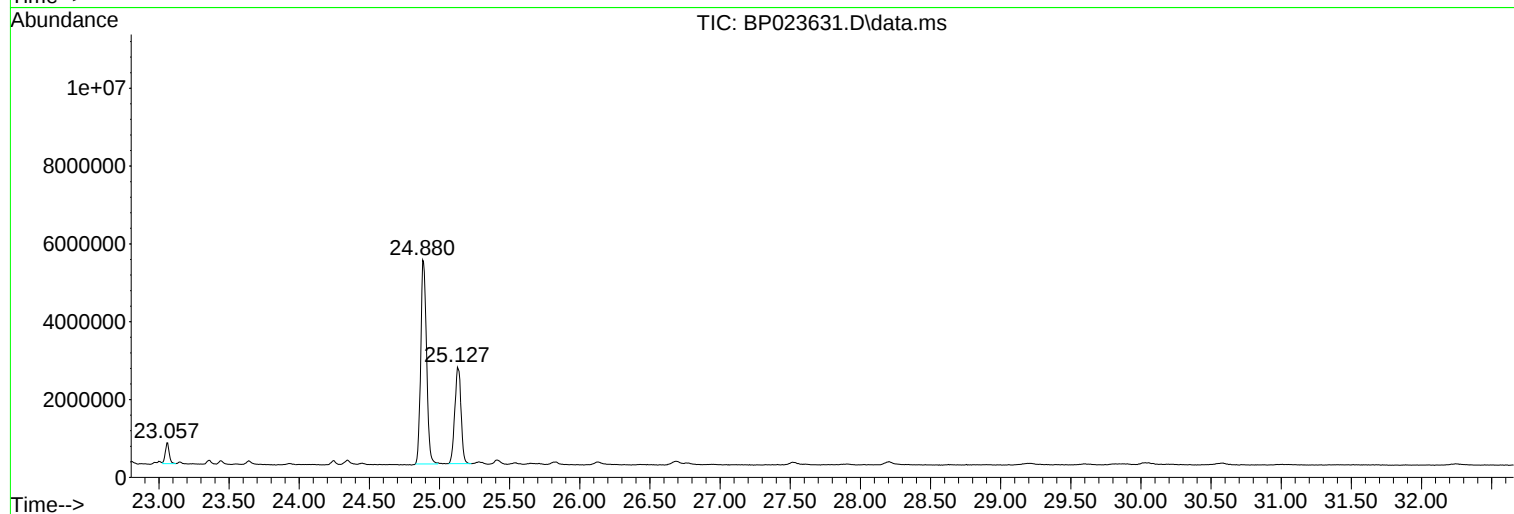
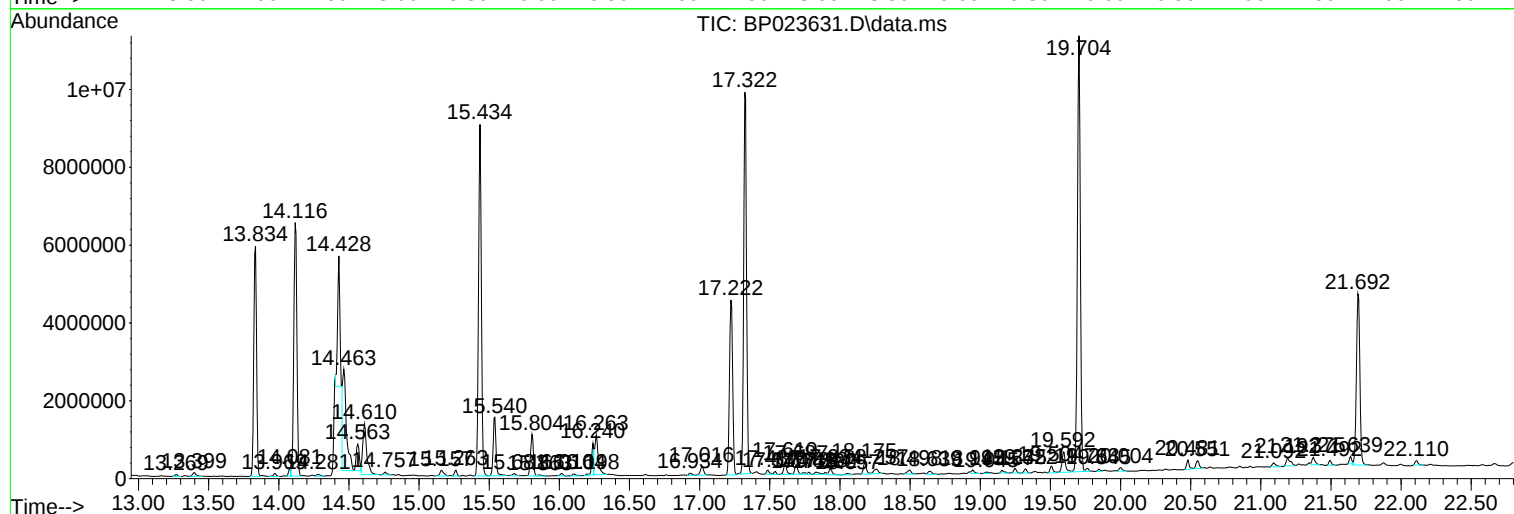
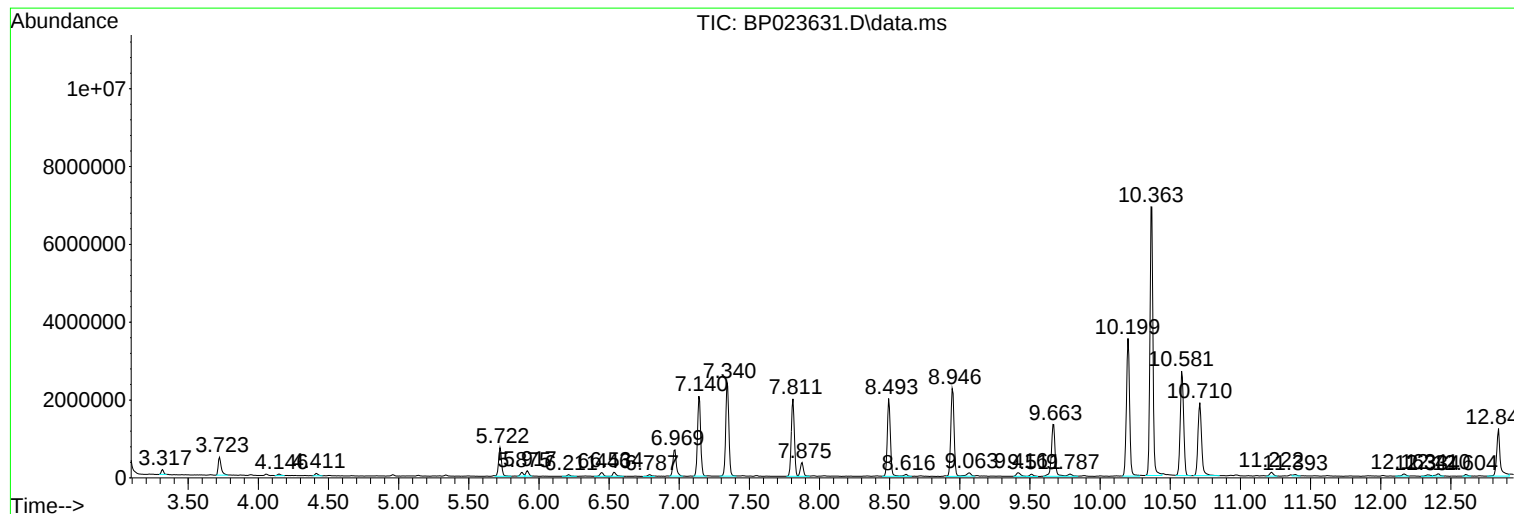
Sum of corrected areas: 177478325

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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP011425.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P



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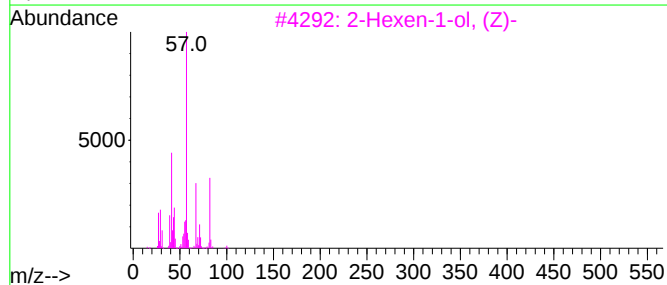
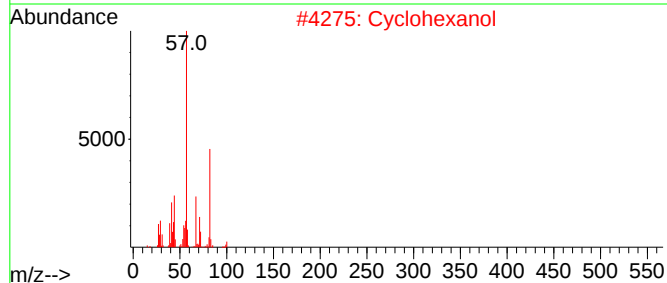
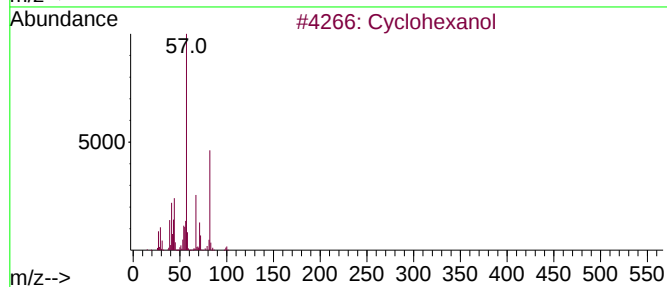
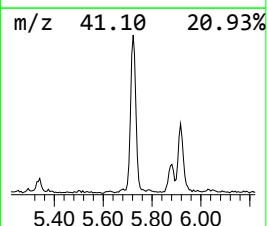
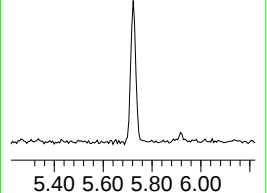
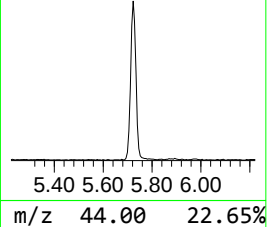
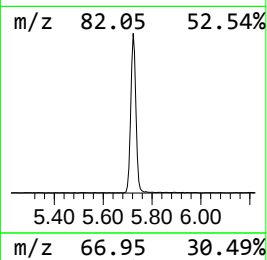
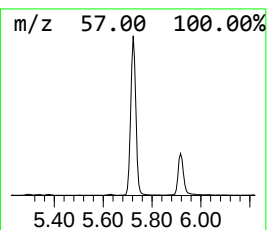
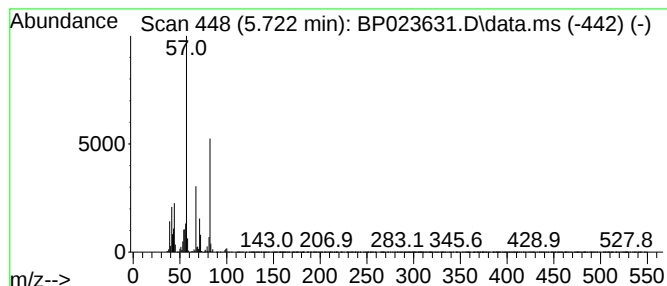
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP011425.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Cyclohexanol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.722	7.23 ng/ul	1097750	1,4-Dichlorobenzene-d4	7.811

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexanol	100	C6H12O	000108-93-0	95
2			Cyclohexanol	100	C6H12O	000108-93-0	91
3			2-Hexen-1-ol, (Z)-	100	C6H12O	000928-94-9	74
4			2-Hexen-1-ol, (E)-	100	C6H12O	000928-95-0	64
5			2-Hexen-1-ol, (E)-	100	C6H12O	000928-95-0	64



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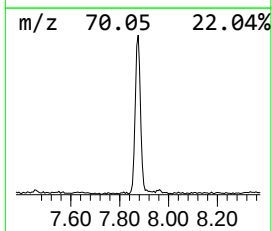
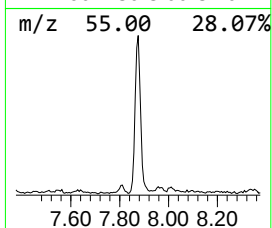
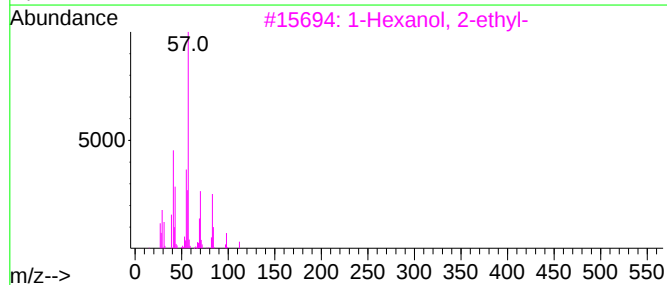
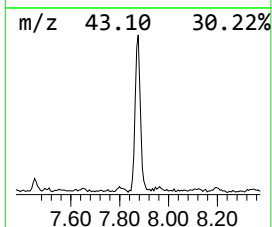
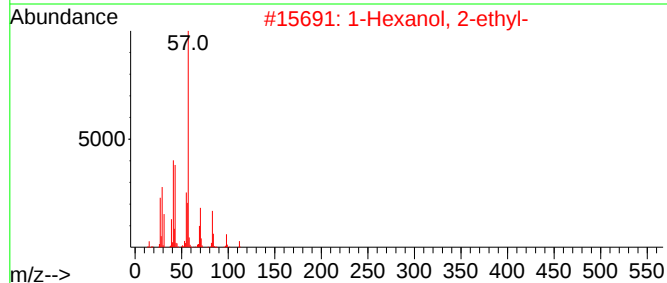
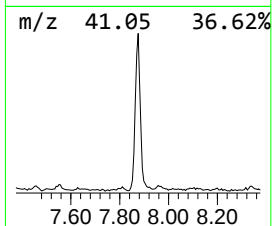
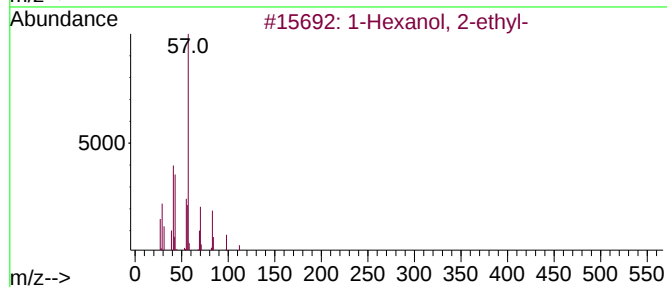
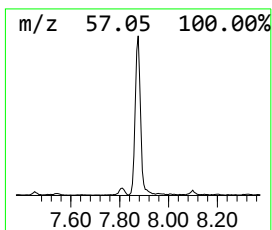
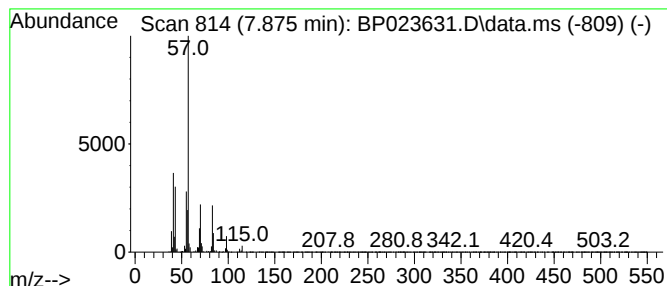
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP011425.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 1-Hexanol, 2-ethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.875	3.56 ng/ul	540925	1,4-Dichlorobenzene-d4	7.811

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	90
2	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	83
3	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	78
4	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	72
5	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	72



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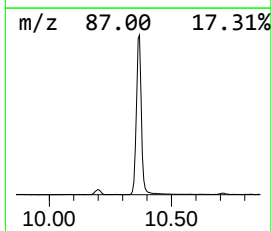
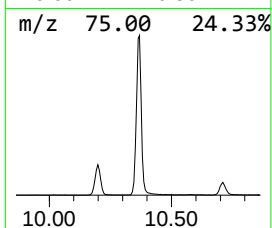
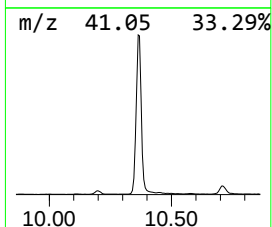
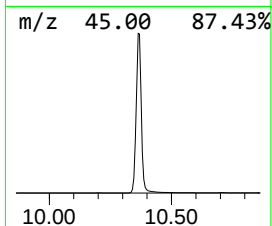
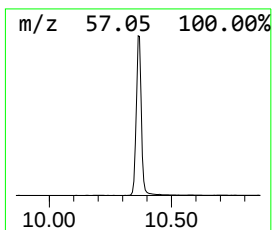
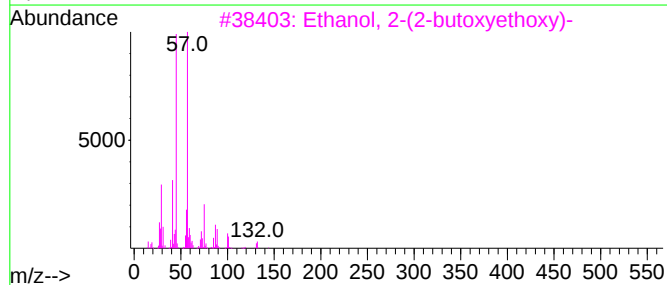
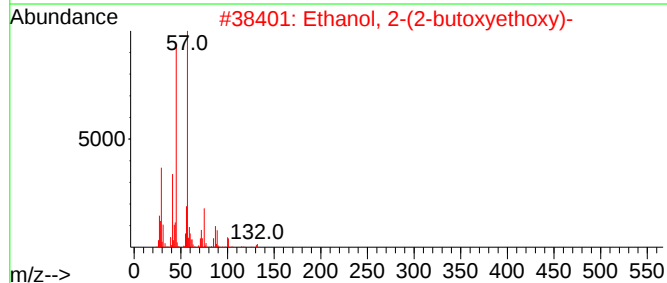
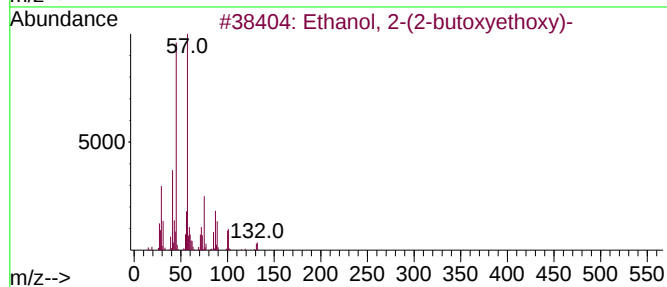
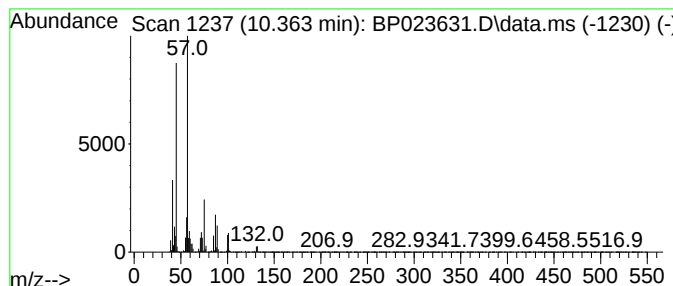
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.363	45.05 ng/ul	10127300	Naphthalene-d8	10.581

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	91
2			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
3			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
4			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	83
5			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011725\
Data File : BP023631.D
Acq On : 17 Jan 2025 16:40
Operator : RC/JU
Sample : Q1086-13
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AF1

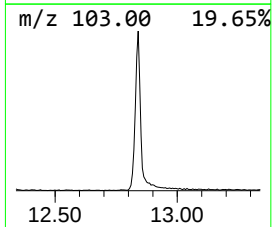
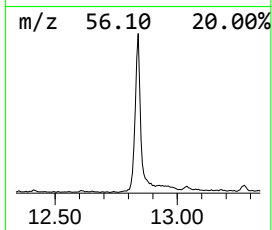
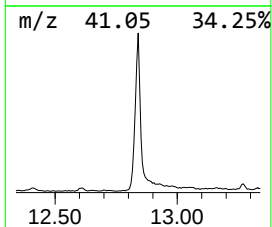
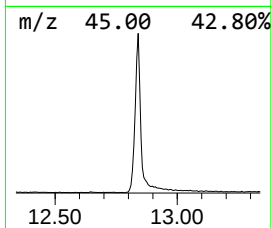
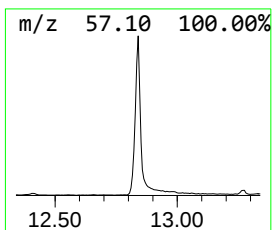
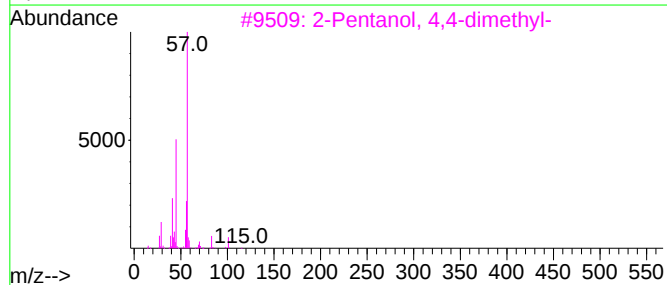
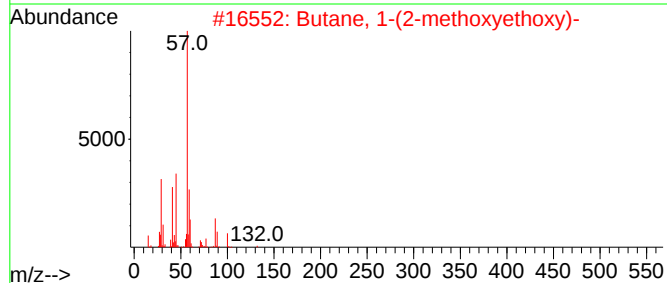
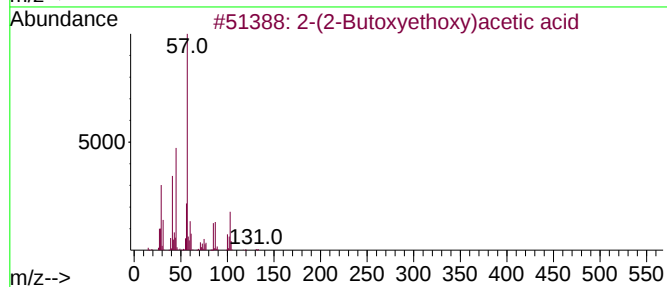
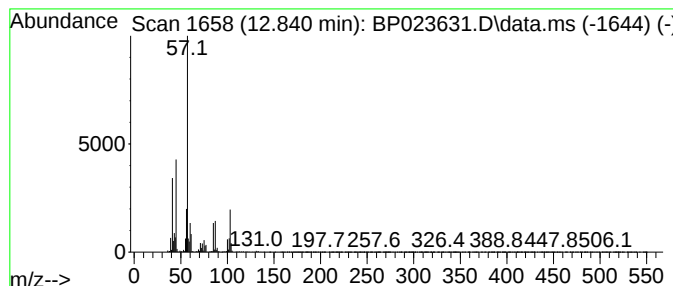
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 2-(2-Butoxyethoxy)acetic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.840	11.31 ng/ul	2072230	Acenaphthene-d10	14.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-(2-Butoxyethoxy)acetic acid	176	C8H16O4	082941-26-2	86
2			Butane, 1-(2-methoxyethoxy)-	132	C7H16O2	013343-98-1	43
3			2-Pentanol, 4,4-dimethyl-	116	C7H16O	006144-93-0	43
4			Butyl 2-(2-butoxyethoxy)acetate	232	C12H24O4	1000366-90-0	42
5			Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011725\
Data File : BP023631.D
Acq On : 17 Jan 2025 16:40
Operator : RC/JU
Sample : Q1086-13
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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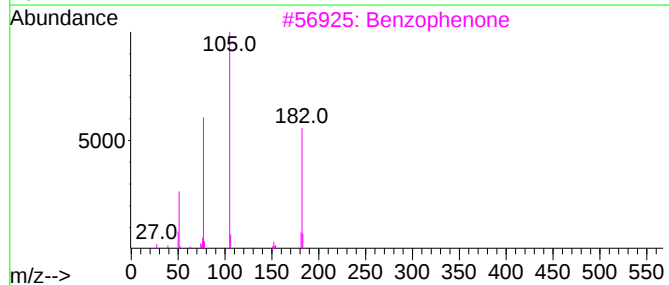
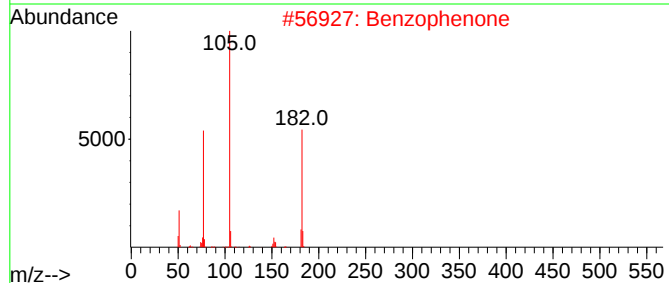
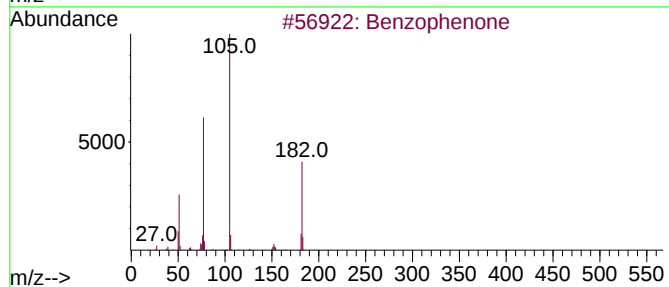
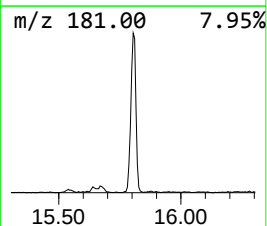
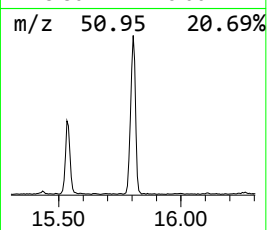
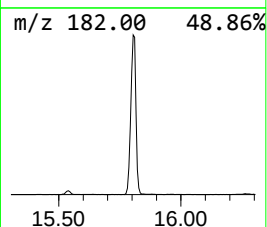
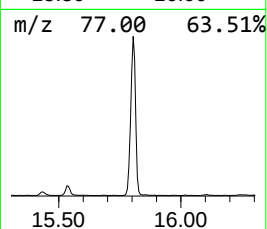
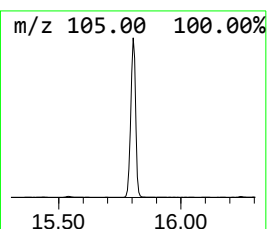
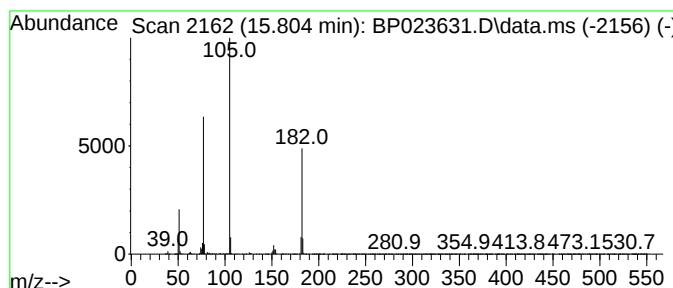
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzophenone Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.804	8.07 ng/ul	1479080	Acenaphthene-d10	14.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	97
2			Benzophenone	182	C13H10O	000119-61-9	96
3			Benzophenone	182	C13H10O	000119-61-9	96
4			Benzophenone	182	C13H10O	000119-61-9	95
5			1,2-Propanedione, 1-phenyl-, 2-o...	163	C9H9NO2	000119-51-7	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011725\
Data File : BP023631.D
Acq On : 17 Jan 2025 16:40
Operator : RC/JU
Sample : Q1086-13
Misc :
ALS Vial : 14 Sample Multiplier: 1

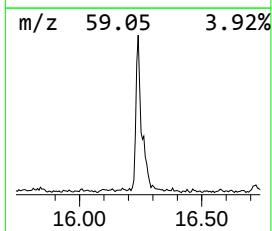
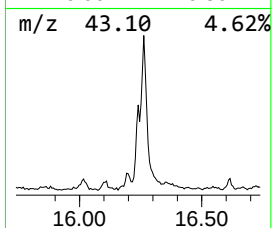
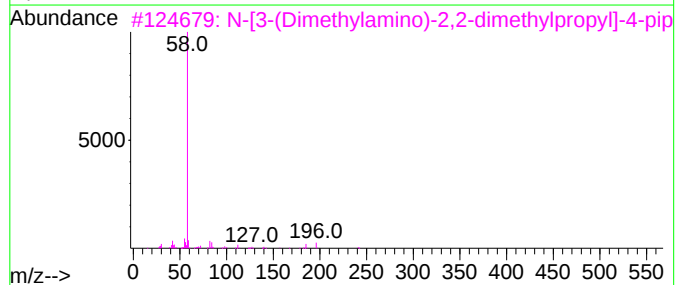
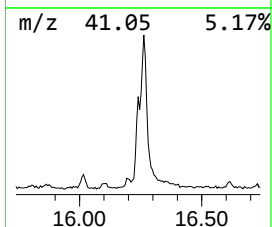
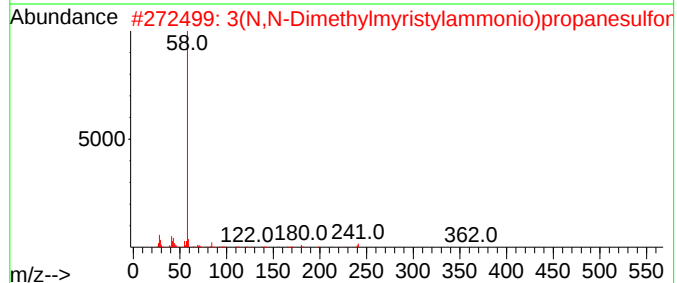
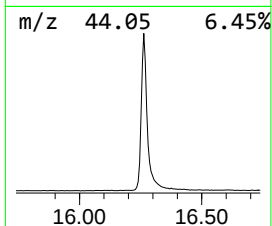
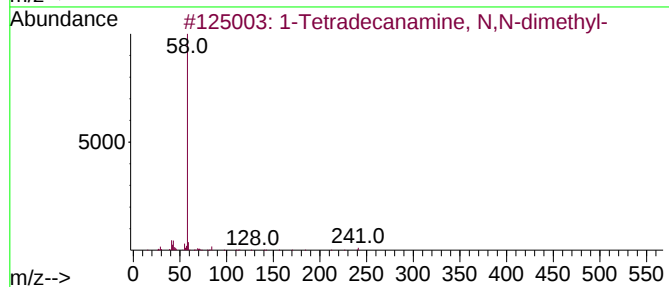
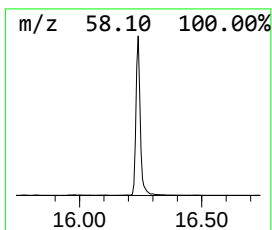
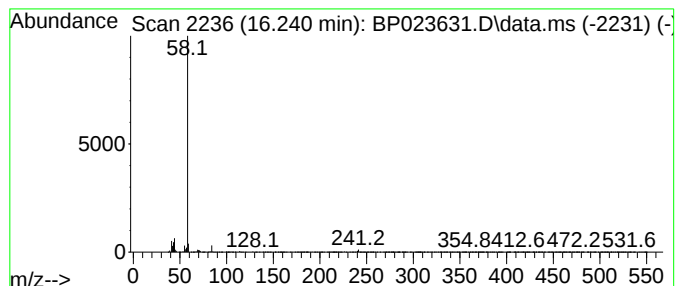
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP011425.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 1-Tetradecanamine, N,N-dime... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.240	2.44 ng/ul	868681	Phenanthrene-d10	17.222	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Tetradecanamine, N,N-dimethyl-	241	C16H35N	000112-75-4	90
2	3(N,N-Dimethylmyristylammonio)pr...	363	C19H41NO3S	014933-09-6	78
3	N-[3-(Dimethylamino)-2,2-dimethyl...	241	C13H27N3O	728932-41-0	78
4	1-Decanaminium, N,N,N-trimethyl-...	235	C13H30ClN	010108-87-9	72
5	1-Tetradecanamine, N,N-dimethyl-	241	C16H35N	000112-75-4	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011725\
Data File : BP023631.D
Acq On : 17 Jan 2025 16:40
Operator : RC/JU
Sample : Q1086-13
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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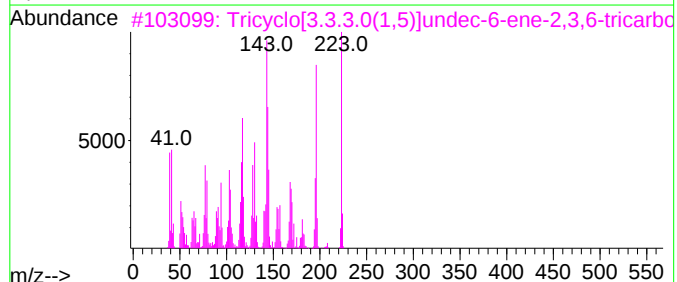
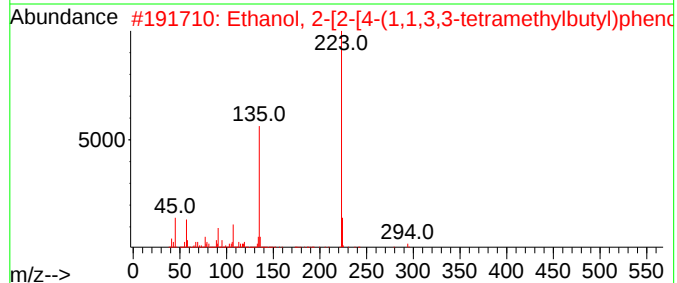
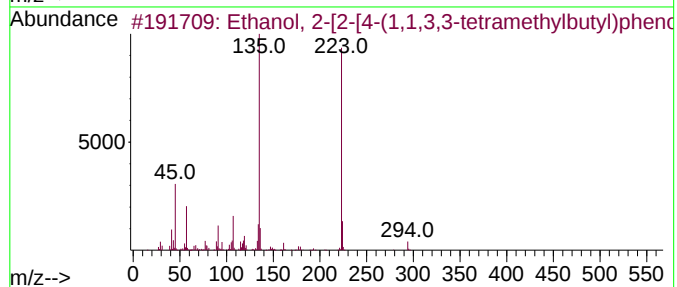
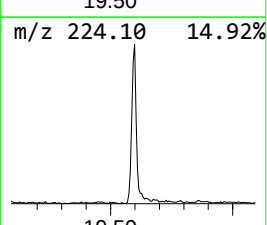
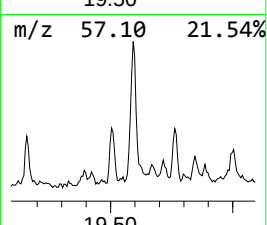
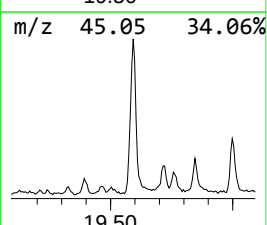
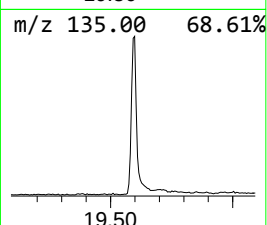
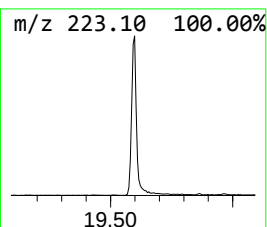
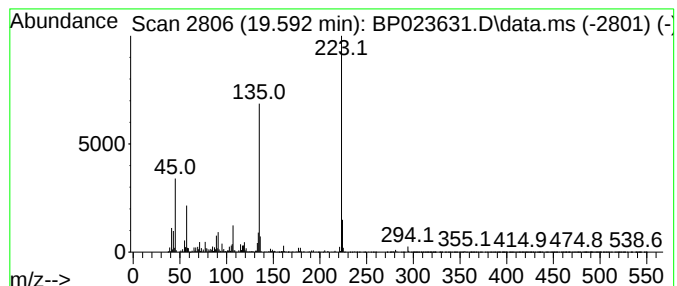
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Ethanol, 2-[2-[4-(1,1,3,3-t... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.592	2.34 ng/ul	880884	Chrysene-d12	21.692

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-[2-[4-(1,1,3,3-tetram...	294	C18H30O3	002315-61-9	97
2			Ethanol, 2-[2-[4-(1,1,3,3-tetram...	294	C18H30O3	002315-61-9	97
3			Tricyclo[3.3.3.0(1,5)]undec-6-en...	223	C14H13N3	118894-71-6	94
4			Ethanol, 2-[2-[4-(1,1,3,3-tetram...	294	C18H30O3	002315-61-9	91
5			Pyrrolo[1,2-a]-1,3,5-triazine-7-...	223	C9H9N3O4	054449-89-7	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011725\
Data File : BP023631.D
Acq On : 17 Jan 2025 16:40
Operator : RC/JU
Sample : Q1086-13
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AF1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP011425.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Cyclohexanol	5.722	7.2	ng/u1	1097750	1	7.811	3036230	20.0
1-Hexanol, 2-et...	7.875	3.6	ng/u1	540925	1	7.811	3036230	20.0
Ethanol, 2-(2-b...	10.363	45.0	ng/u1	10127300	2	10.581	4496460	20.0
2-(2-Butoxyetho...	12.840	11.3	ng/u1	2072230	3	14.428	3664880	20.0
Benzophenone	15.804	8.1	ng/u1	1479080	3	14.428	3664880	20.0
1-Tetradecanami...	16.240	2.4	ng/u1	868681	4	17.222	7107020	20.0
Ethanol, 2-[2-[...	19.592	2.3	ng/u1	880884	5	21.692	7518710	20.0